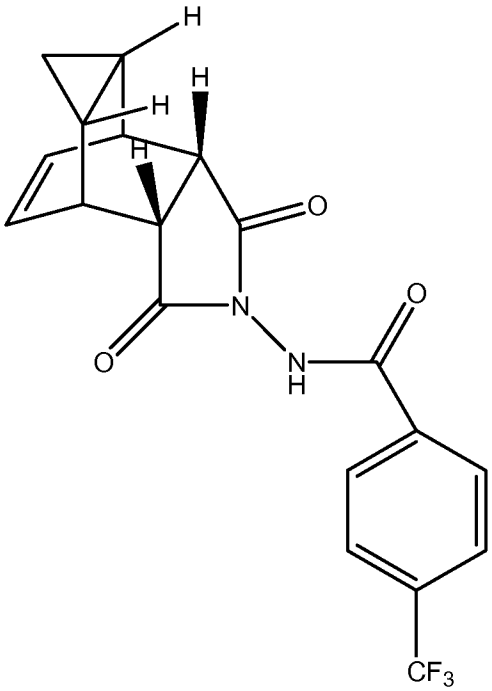
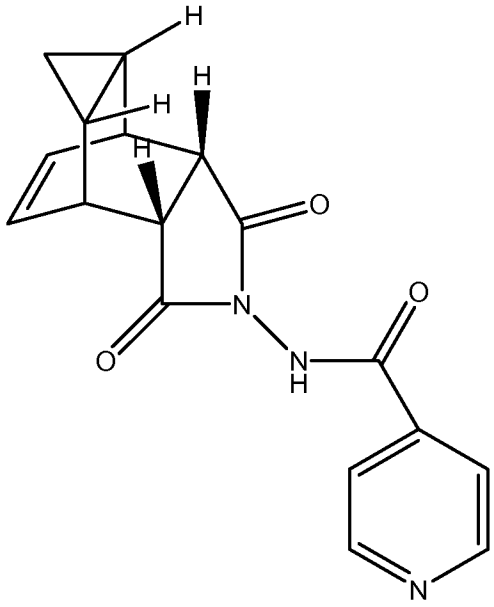
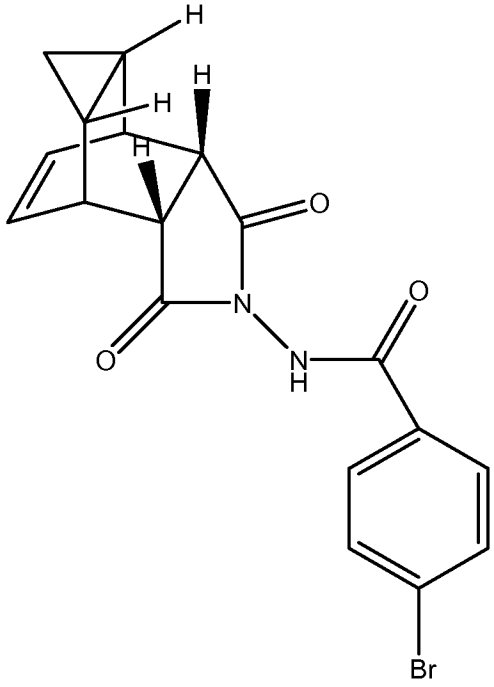
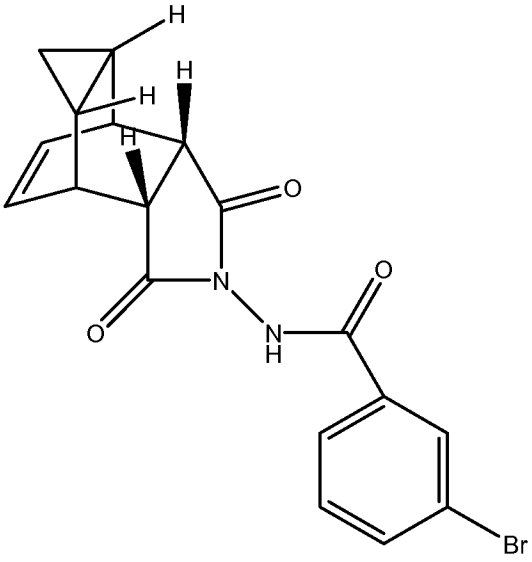
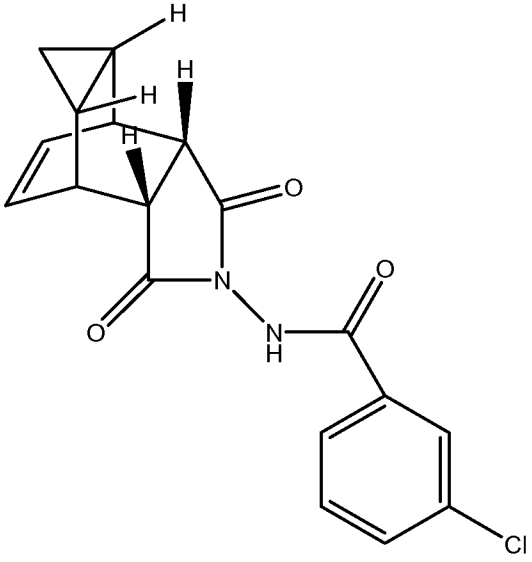
AMENDMENTS TO THE SPECIFICATION

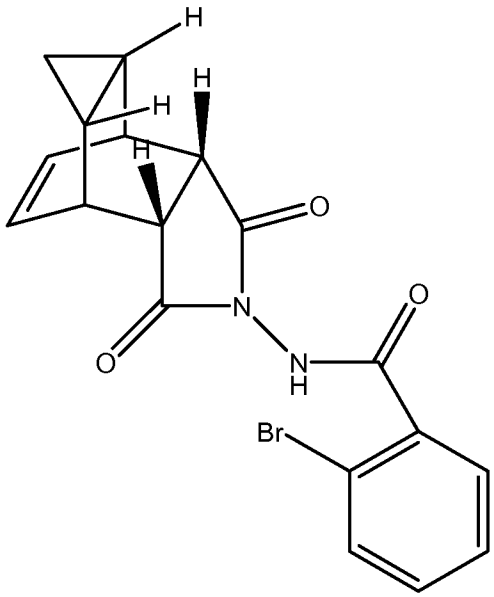
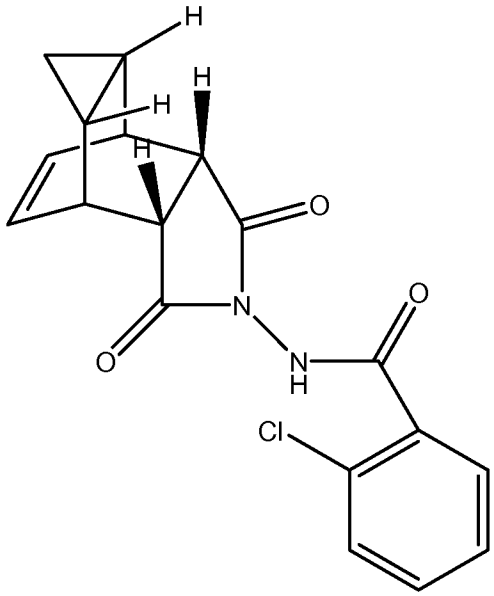
On page 24, after the first paragraph, please amend Table 1 as follows:

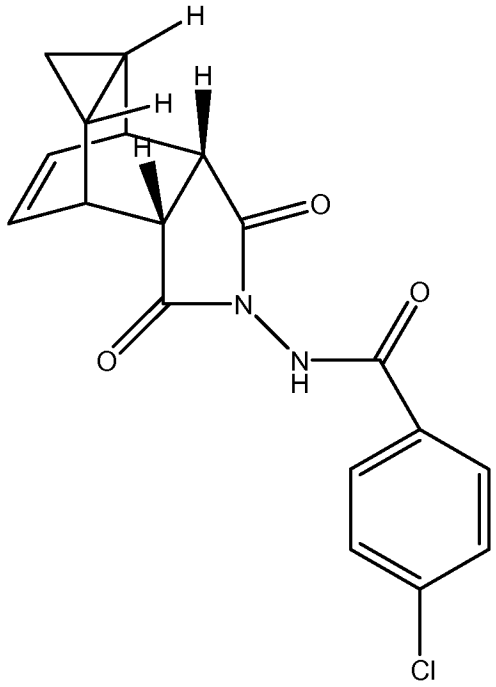
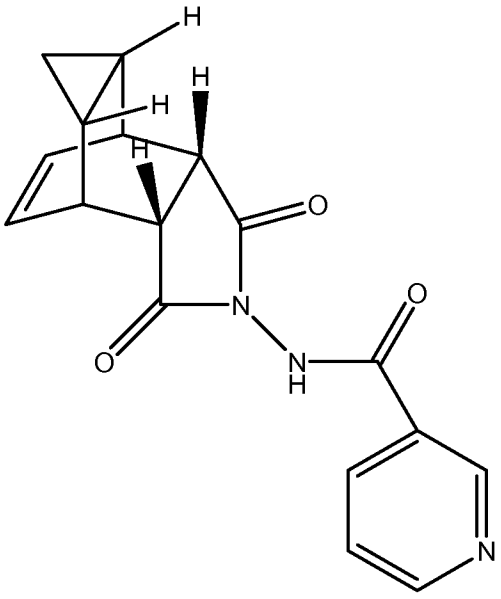
TABLE 1

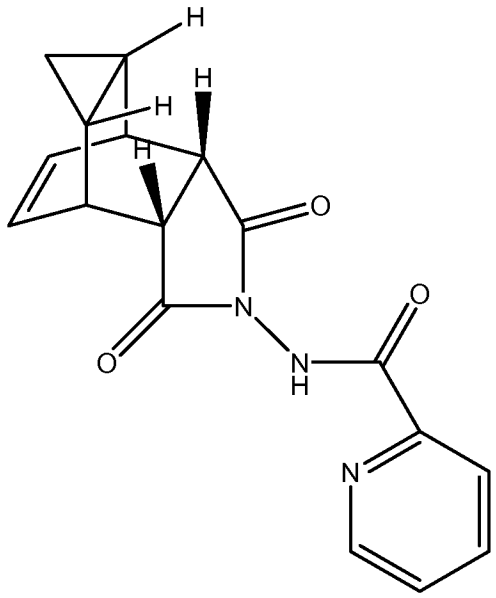
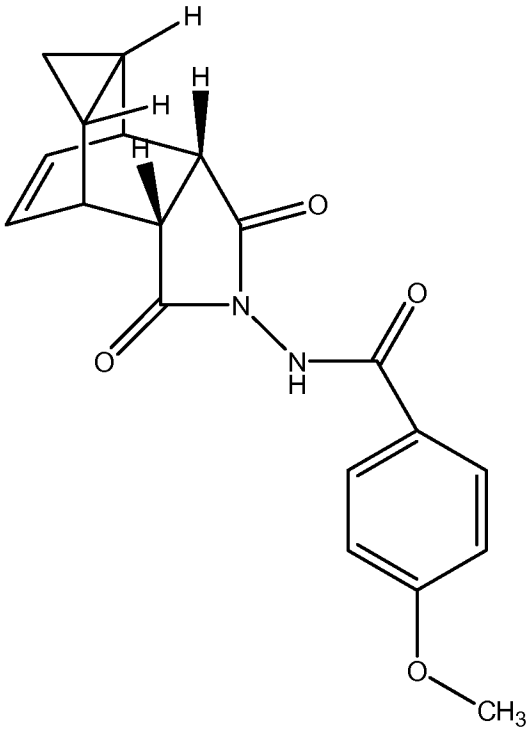
Example Number	R_6	*NMR	**Mass Spec	Name
1		^{1}H NMR in DMSO-d_6: δ 11.35(d, 1H): 11.09(d, 1H); 8.08(d, 2H); 7.92(d, 2H); 5,799(s, 2H); 3.29(brs, 4H), 1.17(m, 2H); 0.26(m, 1H; 0.078(s, 1H)	375 (M-H)-	4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

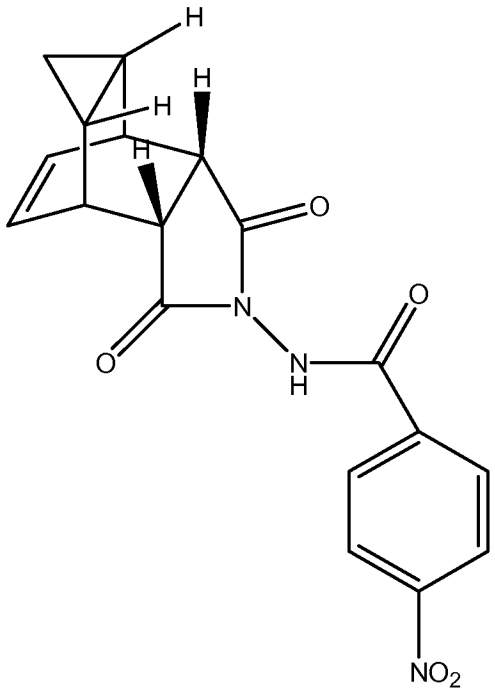
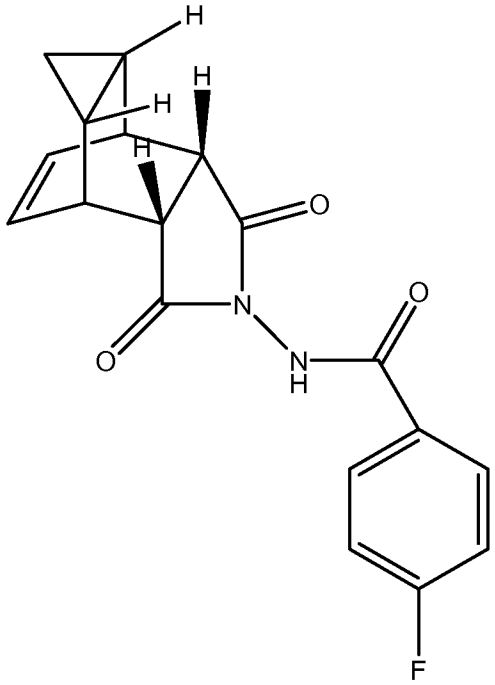
Example Number	R_6	*NMR	**Mass Spec	Name
2		^{1}H NMR in DMSO-d_6: δ 11.41(brs); 11.15(brs); 8.77(d of d, 2H); 7.75(d, 2H); 5,77(brs, 2H); 3.27 (brs, 4H), 1.15(brs, 2H); 0.25(m, 1H; 0.03(brs, 1H)	308 (M-H)-	N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-4-pyridinecarboxamide
3		***	385 (M-H)-	4-Bromo-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

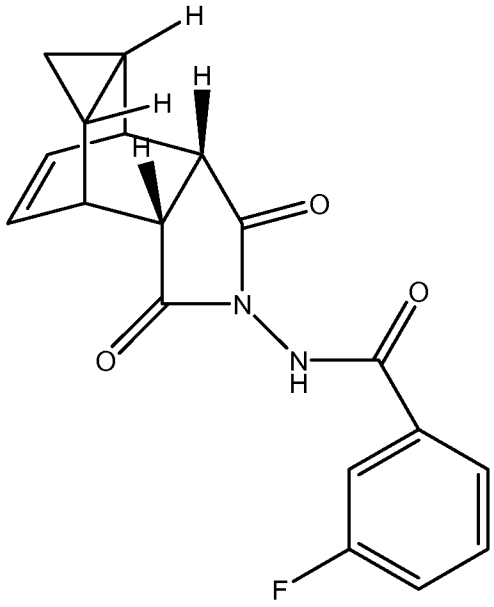
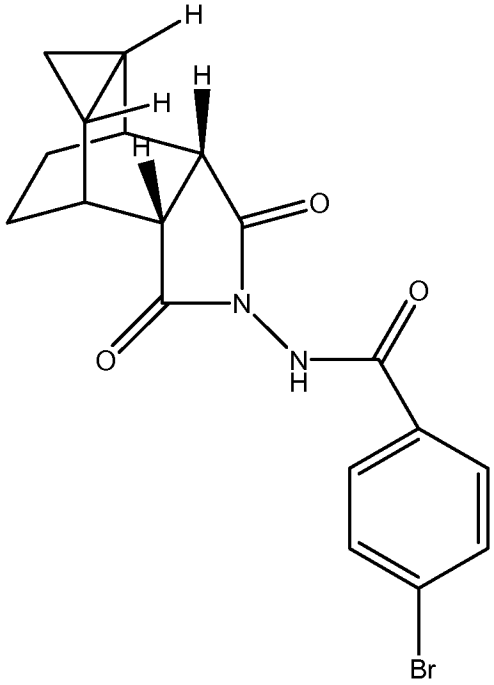
Example Number	R_6	*NMR	**Mass Spec	Name
4		^{1}H NMR in DMSO-d_6: δ 11.13(brd, 1H); 10.89(brd, 1H); 7.99(s, 1H); ~~7.92~~ 7.82 – 7.76(m, 2H); 7.43(t, 1H); 5.72(s, 2H), 3.22-3.08(m, 4H); 1.19(brs, 2H; 0.21(m, 1H); 0.17(brs, 1H)	385 (M-H)-	3-Bromo-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
5		^{1}H NMR in DMSO-d_6: δ 11.21(brd, 1H); 10.98(brd, 1H); 7.92(s, 1H); 7.85(d, 1H); 7.71(d, 1H); 7.58(t, 1H), 5.79(brs, 2H); 3.29-3.15(m, 4H); 1.19-1.15(m, 2H); 0.26(m, 1H); 0.10(brs, 1H)	341 (M-H)-	3-Chloro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

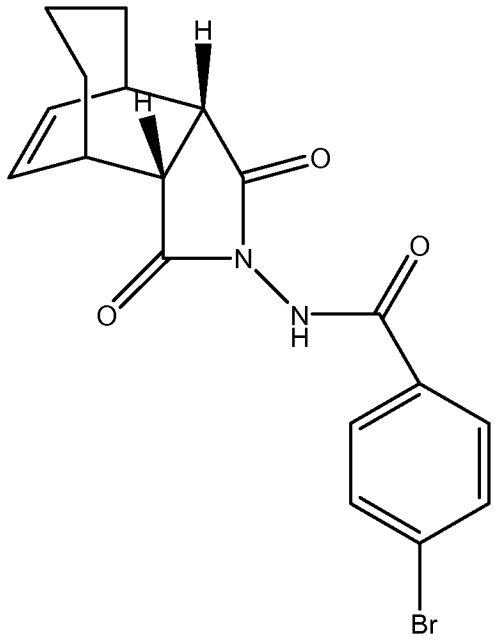
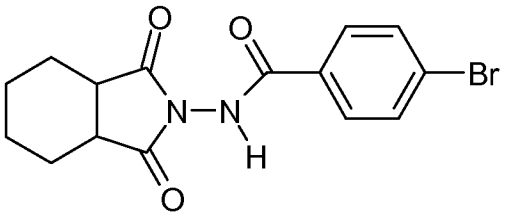
Example Number	R_6	*NMR	**Mass Spec	Name
6		^{1}H NMR in $CDCl_3$: δ 7.74(s, 1H); 7.69(d, 1H); 7.63(d, 1H); 7.41-7.31(m, 2H); 5.84(m, 2H); 3.48(m, 2H), 3.14(s, 2H); 1.19(m, 2H); 0.38-0.20(m, 2H)	385 (M-H-)	2-Bromo-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
7		^{1}H NMR in $CDCl_3$: δ 7.96(s, 1H); 7.83(d, 1H); 7.45(m, 2H); 7.36(m, 1H); 5.86(d, 2H); 3.47(brs, 2H), 3.15(s, 2H); 1.15(brs, 2H); 0.39-0.20(m, 2H)	341 (M-H)-	2-Chloro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

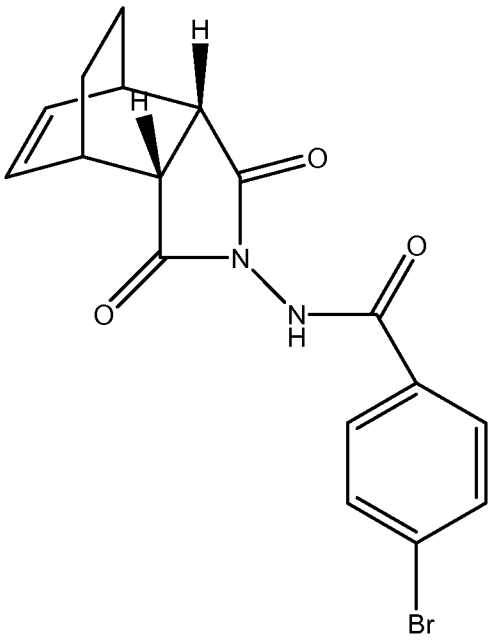
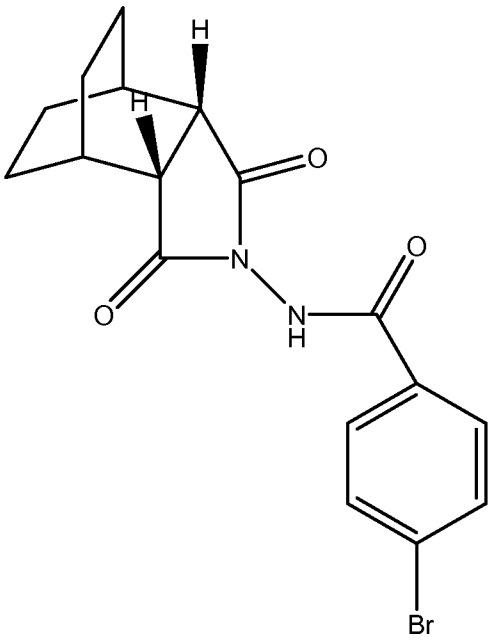
Example Number	R_6	*NMR	**Mass Spec	Name
8		^{1}H NMR in DMSO-d_6: δ 11.16(brd, 1H); 10.91(brd, 1H); 7.90(d, 2H); 7.61(d, 2H); 5.79(s, 2H); 3.28(m, 4H), 1.17(s, 2H); 0.26(m, 1H); 0.07(s, 2H)	341 (M-)H-	4-Chloro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
9		^{1}H NMR in DMSO-d_6: δ 11.33(brd, 1H); 9.04(s, 1H); 8.8(m, 1H); 8.23(d, 1H); 7.56(m, 1H); 5.80(s, 2H), 3.29(m, 4H); 1.17(m, 2H); 0.27(m, 1H); 0.07(s, 1H)	308 (M-)H-	N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-3-pyridinecarboxamide

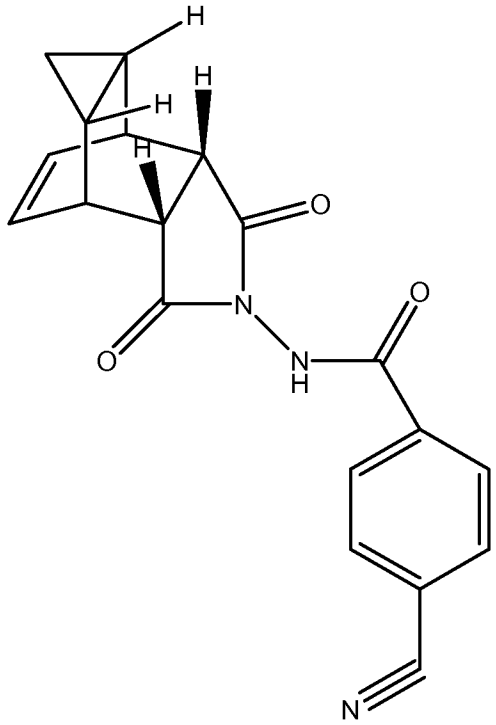
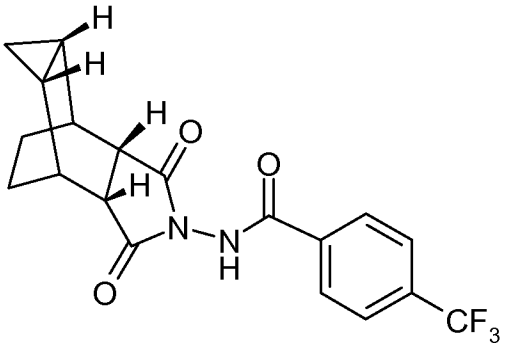
Example Number	R_6	*NMR	**Mass Spec	Name
10		^{1}H NMR in DMSO-d_6: δ 11.11(s, 1H); 8.70(d, 1H); 8.07-8.02(M, 2H); 7.7-7.66(m, 1H); 5.75(m, 2H); 3.295(s, 4H), 1.16(m, 2H); 0.27(m, 1H); 0.10(s, 1H)	308 (M-H)-	N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-2-pyridinecarboxamide
11		^{1}H NMR in DMSO-d_6: δ 10.87(brd, 1H); 7.87(d, 2H); 7.05(d, 2H); 5.78(br, 2H); 3.84(s, 3H); 3.30(s, 4H), 1.16(m, 2H); 0.25(m, 1H); 0.07(brs, 1H)	339 $(M+H)^+$	4-Methoxy-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

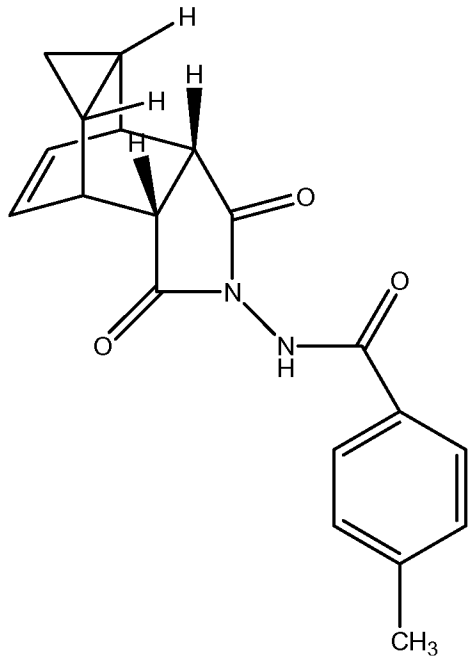
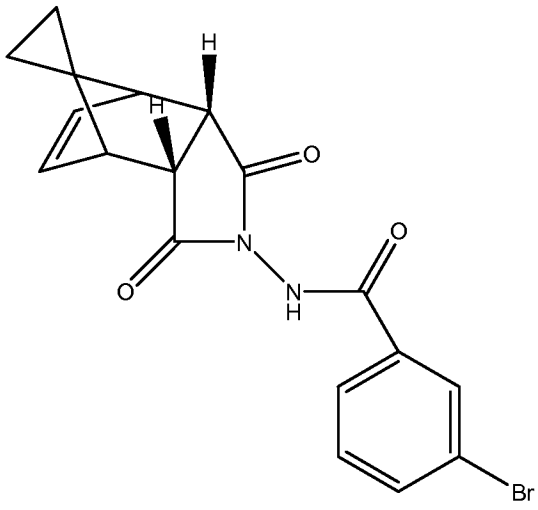
Example Number	R_6	*NMR	**Mass Spec	Name
12		^{1}H NMR in DMSO-d_6: δ 11.537-11.469(brd, 1H); 8.38(d, 2H); 8.12 (d, 2H); 5.80(s, 2H); 3.3(br, 4H); 1.18(s, 2H); 0.27(m, 1H); 0.08(s, 1H)	352 (M-H)-	4-Nitro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
13		^{1}H NMR in DMSO-d_6: δ 11.04(br, 1H); 7.96(s, 2H); 7.367(t, 2H); 5.791(s, 2H); 3.258(4H & H_2O), 1.18(d, 2H); 0.28(m, 1H); 0.09(s, 1H)	327.0 $(M+H)^+$	4-Fluoro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

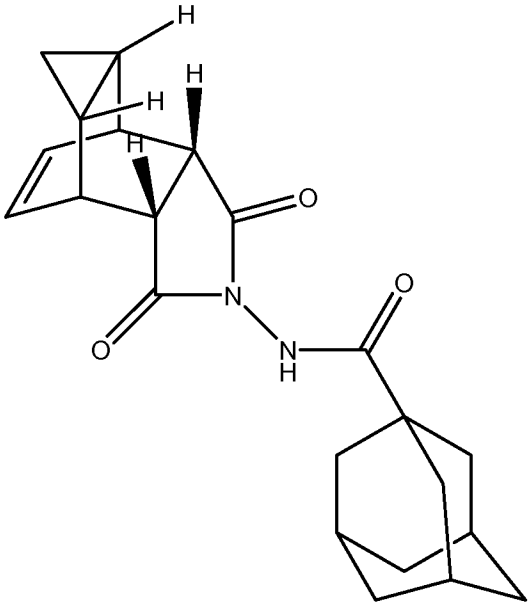
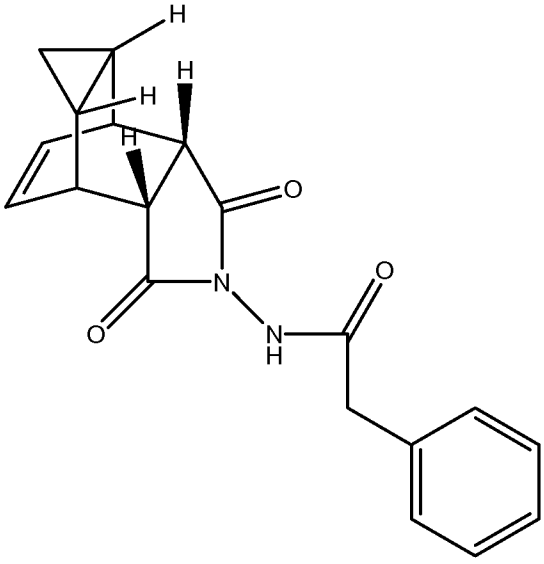
Example Number	R_6	*NMR	**Mass Spec	Name
14		^{1}H NMR in DMSO-d_6: δ 11.176(br, 1H); 7.768-7.459(m, 4H); 5.797(s, 2H); 3.293(H_2O & 4H), 1.174(s, 2H); 0.23(m, 1H); 0.05(s, 1H)	327.0 $(M+H)^+$	3-Fluoro-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
15		***	388.9 (M-H)-	~~4-Bromo-N-(octahydro-1,3-dioxo-4,6-ethanocycloprop[f]isoindol-2(1H)-yl)-benzamide~~ 4-Bromo-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-4,6-ethanocycloprop[f]isoindol-2(1H)-yl)-benzamide

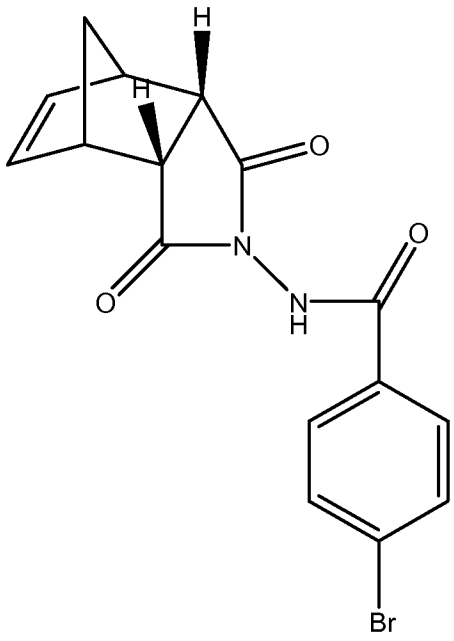
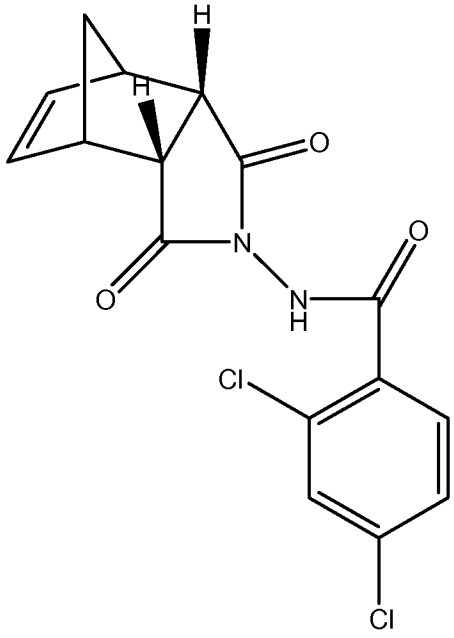
Example Number	R_6	*NMR	**Mass Spec	Name
16		^{1}H NMR in DMSO-d_6: δ 11.14(brd, 1H); 7.85(brd, 2H); 7.76(d, 2H);6.10(brs, 2H) 3.43(brd, 2H), 2.86(brs, 2H); 1.98-1.54(m, 6H)	387 (M-H)-	4-Bromo-N-(1,3-(2H, 3aH)-dioxo-4,8-ethenocyclohepta[c]pyrrolyl)-benzamide
17		^{1}H NMR in DMSO-d_6: δ 11.16(s, 1H); 7.86(d, 2H); 7.78 (d, 2H); 3.14 (brs, 2H); 1.81-1.68 (brm, 4H); 1.42 (br, 4H) ~~5.797(s, 2H); 3.293(H_2O & 4H), 1.174(s, 2H); 0.23(m, 1H); 0.05(s, 1H)~~	350.9 $(M+H)^+$	4-Bromo-N-(octahydro-1,3-dioxo-2H-isoindol-2-yl)-benzamide

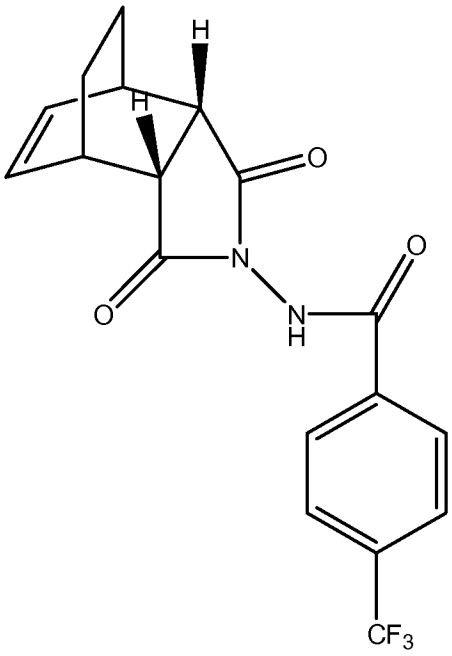
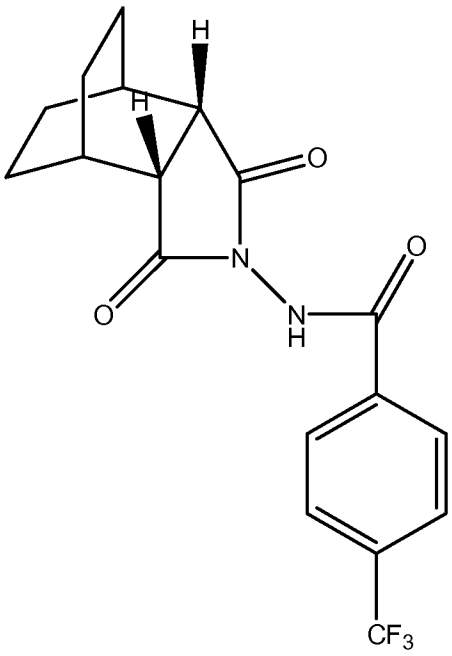
Example Number	R_6	*NMR	**Mass Spec	Name
18		^{1}H NMR in DMSO-d_6: δ 11.05(brd, 1H); 7.83(d, 2H); 7.76(d, 2H); 6.21(s, 2H), 3.15(s, 2H); 3.04(s, 2H);1.66(d, 2H); 1.28(d, 2H)	373 (M-H)-	4-Bromo-N-bicyclo[2.2.2]oct-5-ene-2,3-dicarboximido-benzamide
19		^{1}H NMR in DMSO-d_6: δ 11.15(s, 1H); 7.87(d, 2H); 7.78(d, 2H); 3.07(m, 2H), 2.04(s, 2H); 1.75-1.64(m, 2H); 1.45-1.38(m, 3H)	373 (M-H)-	4-Bromo-N-bicyclo[2.2.2]octane-2,3-dicarboximido-benzamide

Example Number	R_6	*NMR	**Mass Spec	Name
20		^{1}H NMR in DMSO-d_6: δ 11.36(br, 1H); 8.03(s, 4H); 5.79(s, 2H); 3.30(4H + H_2O);2.50(s, 2H); 1.20(s, 2H)	332.1 (M-H)-	4-Cyano-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
21		^{1}H NMR in DMSO-d_6: δ 11.286(br, 1H); 8.13(d, 2H); 8.10(d, 2H); 3.30(4H + H_2O);1.49-1.12 (m, 4H); 0.83(s, 1H); 0.57(s, 1H)	377.0 (M-H)-	4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethanocycloprop[f]isoindol-2(1H)-yl)-benzamide

Example Number	R_6	*NMR	**Mass Spec	Name
22		***	***	4-Methyl-N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide
23		***	***	~~3-Bromo-N-(1',2,2'a,4',7,7'a-hexahydro-1',3'-dioxospiro[cyclopropane-1,8'-[4,7]methano[2H]isoindol]-2'-yl)-benzamide~~ 3-Bromo-N-(1',3,3'a,4',7,7'a-hexahydro-1',3'-dioxospiro[cyclopropane-1,8'-[4,7]methano[2H]isoindol]-2'-yl)-benzamide

Example Number	R_6	*NMR	**Mass Spec	Name
24		***	***	N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-tricyclo[3.3.1.13,7]decane-1-carboxamide
25		***	***	N-(3,3a,4,4a,5,5a,6,6a,-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzeneacetamide

Example Number	R_6	*NMR	**Mass Spec	Name
26		***	***	4-Bromo-N-(1,3,3a,4,7,7a-hexahydro-1,3,-dioxo-4,7-methano-2H-isoindol-2-yl)-benzamide
27		***	***	2,4-Dichloro-N-(1,3,3a,4,7,7a-hexahydro-1,3,-dioxo-4,7-methano-2H-isoindol-2-yl)-benzamide

Example Number	R_6	*NMR	**Mass Spec	Name
28		^{1}H NMR in DMSO-d_6: δ 11.37(br, 1H); 8.10(d, 2H); 7.94 (d, 2H); 6.22(s,2H);3.17 (s, 2H); 3.05(s, 2H); 1.66(m, 2H); 1.29 (m, 2H)	365.0 $(M+H)^+$	4-Trifluoromethyl-N-bicyclo [2.2.2]oct-5-ene-2,3-dicarboximido-benzamide
29		^{1}H NMR in DMSO-d_6: δ 11.33(s, 1H); 8.14(d, 2H);8.11 (d, 2H);3.29 (s,4H); 2.05(s, 2H); 1.76-1.65(m, 4H); 1.42 (s, 2H)	367.0 $(M+H)^+$	4-Trifluoromethyl-N-bicyclo [2.2.2]octane-2,3-dicarboximido-benzamide

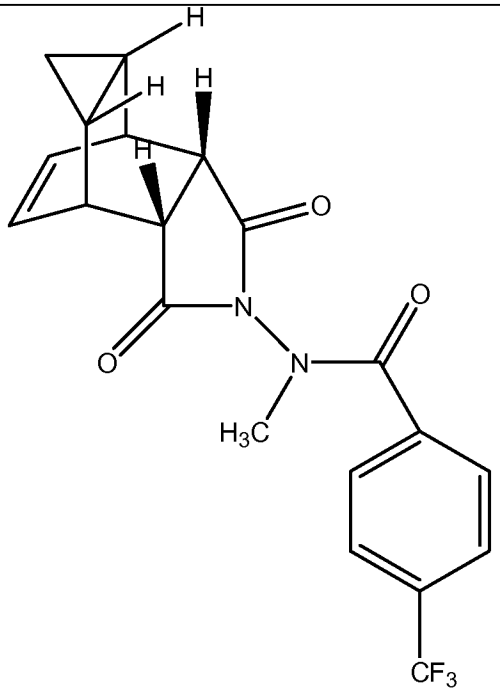
* All 1H NMR and 13C NMR spectra were acquired on a Varian Mercury VX300 Spectrometer and referenced to tetramethysilane (TMS) unless indicated otherwise. Chemical shifts and coupling constants are reported in parts per million (ppm) and Hertz (Hz), respectively, Multiplicities indicated are: s=singlet, d=doublet, t=triplet, q=quartet, m=multiplet, dd=doublet of doublets, and br indicates a broad signal.

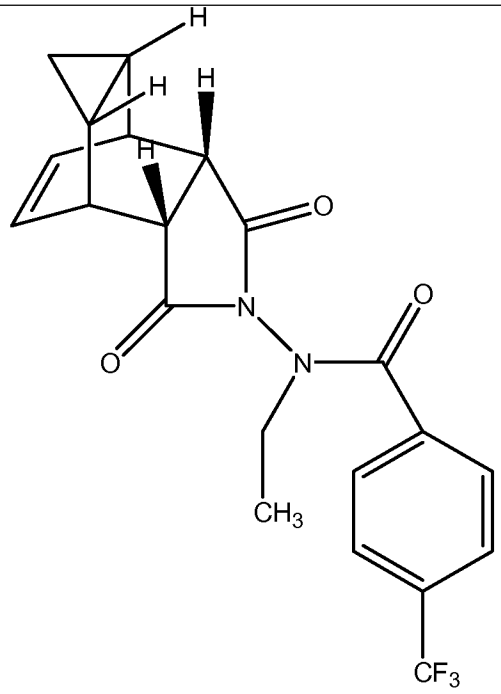
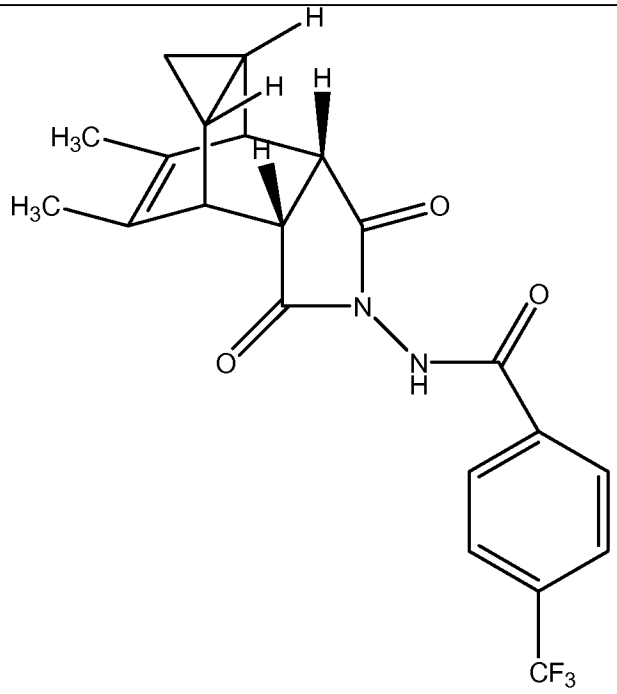
**Mass Spectroscopy data is expressed as a mass to charge ratio (m/z) for either (M+1) or (M-1) molecular ion.

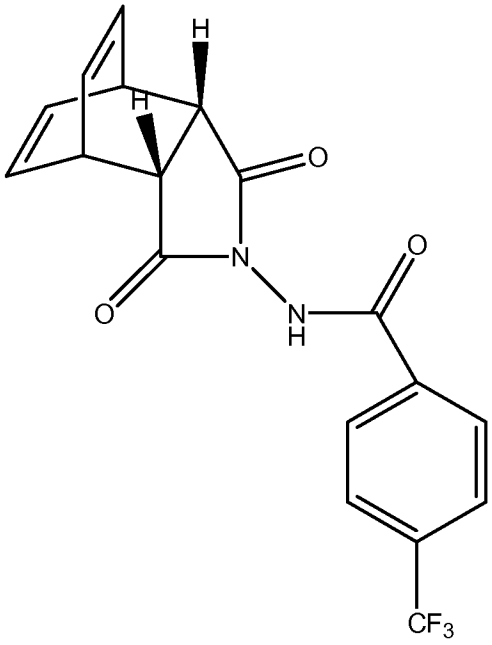
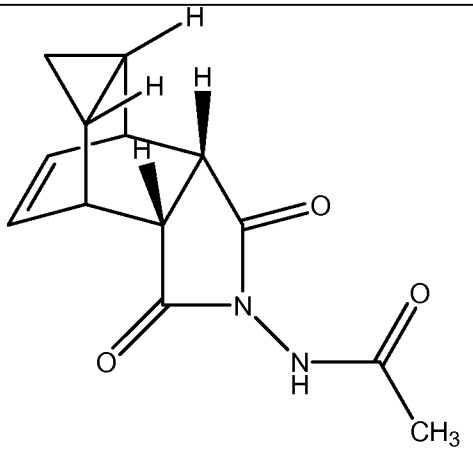
***indicates that data were not collected.

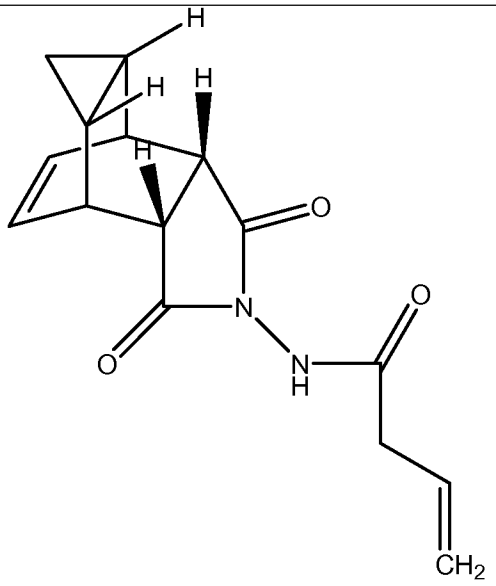
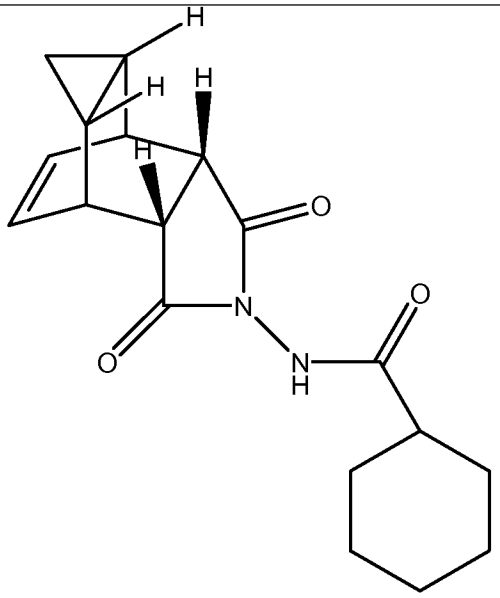
On page 32, please amend Table 2 as follows:

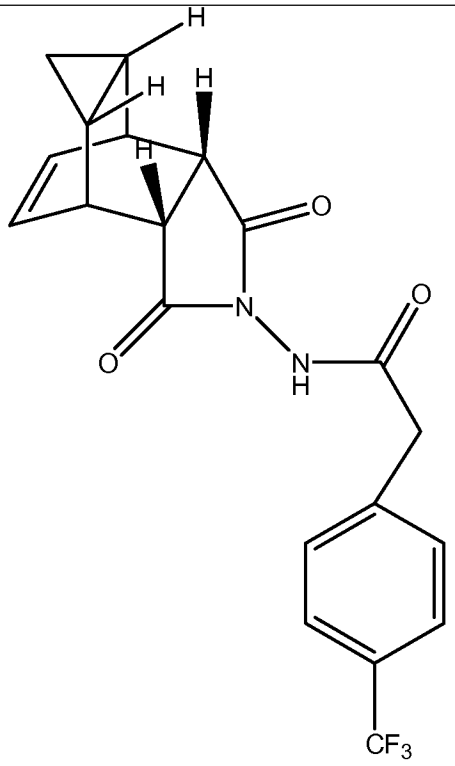
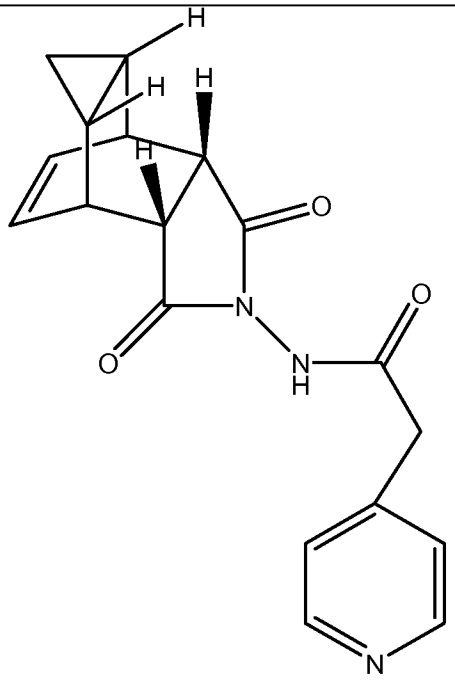
TABLE 2

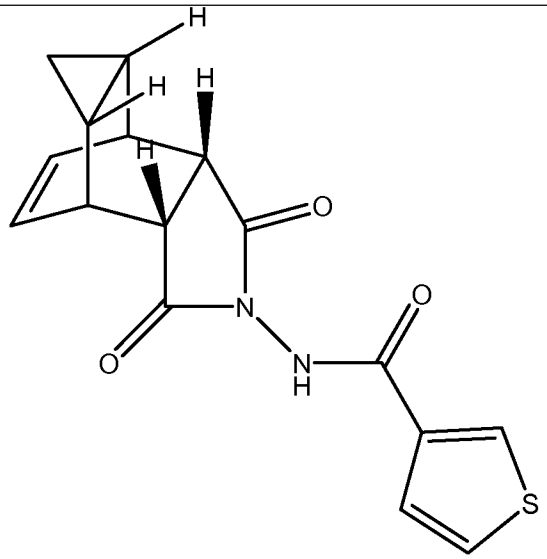
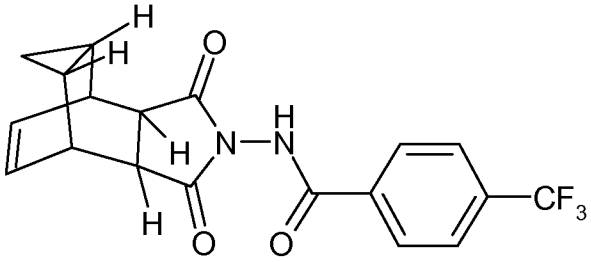
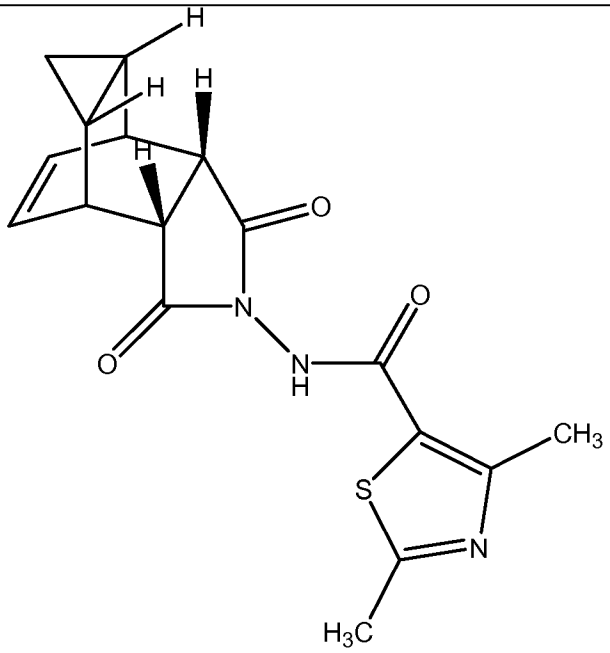
Example Number	Structure	Name
30		4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-N-methylbenzamide

Example Number	Structure	Name
31		4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-N-ethylbenzamide
32		4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzamide

Example Number	Structure	Name
33		4-Trifluoromethyl-N-(3a,4,7,7a-tetrahydro-4,7-etheno-1H-isoindol-2(1H)-yl)-benzamide
34		N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-acetamide

Example Number	Structure	Name
35		N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-but-3-enamide
36		N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-cyclohexanecarboxamide

Example Number	Structure	Name
37		4-Trifluoromethyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-benzylacetamide
38		4-Pyridyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-acetamide

Example Number	Structure	Name
39		~~3-Thienyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-N-methylbenzamide~~ 3-Thienyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-7,8-dimethyl-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)acetamide
40		4-(Trifluoromethyl)-N-[(3aR,4S,4aS,5aR,6R,6aS)-3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl]-benzamide
41		2,4-Dimethyl-N-(3,3a,4,4a,5,5a,6,6a-octahydro-1,3-dioxo-4,6-ethenocycloprop[f]isoindol-2(1H)-yl)-thiazole-5-carboxamide